

ACHIEVING CONSENSUS IN SELF-ORGANIZING WIRELESS SENSOR NETWORKS: THE IMPACT OF NETWORK TOPOLOGY ON ENERGY CONSUMPTION

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ABSTRACT

Achieving consensus on common global parameters through totally decentralized algorithms is a topic that has attracted considerable attention in the last few years. Several algorithms have been developed, among which the most popular is the average consensus method. The main advantage of these approaches is that they do not require a fusion center. But, on the other hand, they are typically based on iterative algorithms, whose energy consumption is proportional to the time necessary to achieve consensus. This time depends on the network topology, as well as on the transmit power of each node. In this paper, we show that there exists an optimal transmit power that minimizes the overall energy consumption necessary to achieve the global estimate within a given accuracy and that this power depends on the network topology.

Index Terms— Average consensus, energy consumption, algebraic connectivity, distributed estimation.

1. INTRODUCTION

Achieving consensus on a common global parameter through totally distributed algorithms is a key problem in wireless sensor networks. In centralized schemes, the need to send sensor data to a fusion center causes congestion around the sink node. With distributed schemes, this congestion is not created, and the network becomes more resilient to node failures and attacks (e.g., an attack on the fusion node could be catastrophic). Distributed agreement algorithms have been studied for a long time; see, e.g. [1]. In recent years, they have received considerable attention, in view of their potential application in sensor networks. The so called *consensus average* method is an example of an algorithm achieving consensus in a totally distributed way [2]. Significant contribution have also come from the context of multiagent coordination and flocking [3].

One of the most critical aspects of these consensus algorithms is that they are iterative algorithms where, at each

step, the network nodes exchange data among each other to achieve agreement. The network need not be fully connected; indeed, consensus is achieved locally, with local islands of agreement expanding, and resulting in global agreement, under appropriate conditions. Consider a network of N nodes. Let p_{Ti} denote the transmit power of the i -th node, assumed to be constant across the time necessary to achieve consensus. Then, the overall energy spent to reach a common estimate (decision), within a given accuracy (reliability), is $\mathcal{E} = T_{conv} \sum_{i=1}^n p_{Ti}$, where T_{conv} is the time necessary to achieve global consensus. The convergence rate has been derived under a variety of situations and it is strictly related to the network topology. More specifically, modeling the network as a graph described by the Laplacian \mathbf{L} , the convergence properties of distributed consensus algorithms depend on the graph spectral properties, i.e., the set of eigenvalues $\lambda_i(\mathbf{L})$, $i = 1, \dots, n$, of \mathbf{L} . In particular, if the network is connected, the smallest eigenvalue $\lambda_1(\mathbf{L})$ is zero and it has algebraic multiplicity one. Furthermore, if the network is connected, the second smallest eigenvalue $\lambda_2(\mathbf{L})$, known as the network *algebraic connectivity*, provides important properties about network connectivity. More specifically, in most consensus algorithms, as we will see in the next section, the convergence rate is directly proportional to $\lambda_2(\mathbf{L})$. Given the critical role played by $\lambda_2(\mathbf{L})$, many papers have concentrated on how to maximize $\lambda_2(\mathbf{L})$ in order to minimize the convergence time, either *rewiring* the networks, i.e. by changing the network topology, as in [4], or by assigning different weights to each link and then optimizing the weight distribution, as in [5].

In general, however, the most critical parameter in wireless sensor networks is typically the energy consumption, which is directly proportional to the convergence time and to the transmit power. On one hand, to save energy, we would like to use the minimum transmit power that ensures network connectivity. But a small transmit power has an effect on the network topology, as it leads to a reduced number of links and, as a consequence, to a small algebraic connectivity $\lambda_2(\mathbf{L})$. Hence, a small individual transmit power implies a long convergence time. Conversely, to reduce the convergence time, the network should have a high connectivity, but this requires a large transmit power. It is then intuitive to expect an optimal trade-off. For simplicity, here we assume that all the

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nodes have the same transmit power. The goal of this paper is to derive the transmit power that yields the minimum energy consumption, for a given topology. We will consider networks with deterministic or random spatial distribution of the nodes. Interestingly, the comparison will clarify whether it is better to distribute sensors randomly or over a regular grid, for any given density and coverage area.

2. CONSENSUS ALGORITHMS

Let us denote with \mathbf{x}_0 the $n \times 1$ vector containing the data collected by all the nodes. Consensus algorithms typically work by endowing each node with a dynamical system that evolves in time as a function of its own state and of a combination of the states of the other nodes. Each system is initialized with the local measurement. If we denote with $\mathbf{x}(t)$ the state vector, at time t , the evolution of all the states is described by the following equation [4]

$$\dot{\mathbf{x}}(t) = -\mathbf{L} \mathbf{x}(t) \quad (1)$$

where \mathbf{L} is the Laplacian of the graph. The Laplacian is defined as follows. Given an oriented graph \mathcal{G}^1 composed of N vertices and E edges, the incidence matrix \mathbf{B} is the $N \times E$ matrix such that $[\mathbf{B}]_{ij} = 1$ if the edge j is incoming to vertex i , $[\mathbf{B}]_{ij} = -1$ if the edge j is outgoing from vertex i , and 0 otherwise. Given the $N \times 1$ vector $\mathbf{1}$, composed of all ones, it is easy to check that the incidence matrix satisfies $\mathbf{1}^T \mathbf{B} = \mathbf{0}^T$. Given \mathbf{B} , the symmetric $N \times N$ matrix $\mathbf{L} \triangleq \mathbf{B} \mathbf{B}^T$, is called the *Laplacian* of \mathcal{G} , and it is independent of the graph orientation. If we associate a positive number a_i to the i -th edge and we build the diagonal matrix $\mathbf{D}_a \triangleq \text{diag}(\mathbf{a})$, with $\mathbf{a} \triangleq [a_1, \dots, a_E]^T$, the Laplacian of the corresponding graph (called the *weighted Laplacian*) is written as $\mathbf{L}_a \triangleq \mathbf{B} \mathbf{D}_a \mathbf{B}^T$. The Laplacian and the weighted Laplacian have several important properties, such as: \mathbf{L} is always positive semi-definite with the smallest eigenvalue equal to 0; the algebraic multiplicity of the null eigenvalue is equal to the number n_c of connected components of the graph. Hence, if the graph is connected, $n_c = 1$ and $\text{rank}(\mathbf{L}) = N - 1$.

By exploiting the properties of the Laplacian, it is easy to check that if the state vector $\mathbf{x}(t)$ is initialized with the local measurement, i.e. $\mathbf{x}(0) = \mathbf{x}_0$, and the network is connected, then $\mathbf{x}(t)$ converges to the average consensus vector

$$\mathbf{x}(t) \rightarrow \frac{1}{N} \mathbf{1} \mathbf{1}^T \mathbf{x}_0. \quad (2)$$

In words, all states converge to the average value.

In the presence of coupling noise, i.e. adding a noise vector $\mathbf{v}(t)$ to (1), the system still converges in the mean. However, as already noticed in [5], the running average $a(t) :=$

$1/N \sum_{i=1}^N x_i(t)$ is a random walk in this case, so that its variance increases unboundedly with time. To alleviate this problem, an algorithm was derived in [5] to compute the weights a_i that minimize the mean square deviation, i.e. $\Delta(t) := E\{\sum_{i=1}^n (x_i(t) - a(t))^2\}$.

To avoid the random walk problem, it is sufficient to associate the average consensus to the derivative of the state, rather than to the state itself. An example of this alternative approach was proposed in [6], [7], where the state evolution was described by the following equation, for $i = 1, \dots, N$,

$$\dot{x}_i(t) = g_i(x_{i0}) + \frac{K}{c_i} \sum_{j=1}^N a_{ij} f(x_j(t) - x_i(t)), \quad (3)$$

where $g_i(x_{i0})$ is a function of the local measurement x_{i0} , K is a global control loop gain; c_i is a local coefficient that quantifies the attitude of the i -th sensor to adapt its values as a function of the signals received from the other nodes: The larger c_i is, the less likely is the i -th node to change its original decision $g_i(x_{i0})$. The function $f(\cdot)$ is, in general, a nonlinear function describing the mutual coupling between sensors and it depends on the radio interface². The running decision, or estimate, of each sensor is encoded in its state derivative $\dot{x}_i(t)$.

Using the incidence matrix \mathbf{B} , we can rewrite the system (3) in compact form as

$$\dot{\mathbf{x}}(t) = \mathbf{g}(\mathbf{x}_0) - K \mathbf{D}_c^{-1} \mathbf{B} \mathbf{D}_A f(\mathbf{B}^T \mathbf{x}(t)), \quad (4)$$

where $\mathbf{x}(t) \triangleq [x_1(t), \dots, x_N(t)]^T$, $\mathbf{D}_c \triangleq \text{diag}\{c_1, \dots, c_N\}$; \mathbf{D}_A is an $E \times E$ diagonal matrix, whose diagonal entries are all the weights a_{ij} , indexed from 1 to E ; the symbol $f(\mathbf{x})$ has to be interpreted as the vector whose k -th component is $f(x_k)$. It was proved in [7] that, if the network is connected, the function $f(x)$ is monotonically increasing and odd³, and K is greater than the threshold

$$K_U = \frac{2 \|\mathbf{D}_c \Delta \omega\|_2}{f_{\max} \lambda_2(\mathbf{L}_a)}, \quad (5)$$

where $f_{\max} = \max f(x)$, $\Delta \omega := \mathbf{g}(\mathbf{x}_0) - \omega^* \mathbf{1}$, with

$$\omega^* = \frac{\mathbf{c}^T \mathbf{g}(\mathbf{x}_0)}{\mathbf{1}_N^T \mathbf{c}} = \frac{\sum_{i=1}^N c_i g_i(x_{0i})}{\sum_{i=1}^N c_i}, \quad (6)$$

then the networks achieves global agreement in the sense that all the state derivatives converge to the common value ω^* , i.e.

$$\dot{x}_i(t) \rightarrow \omega^*, \quad \forall i. \quad (7)$$

This result was used in [6] and [8] to obtain globally optimal decision schemes, for estimation or detection, respectively,

²Without loss of generality, $f(x)$ is normalized so that $df(0)/dx = 1$. A different value of $df(0)/dx$ can always be included in K .

³Linear coupling, i.e., $f(x) = x$, is then included as a particular case.

¹An orientation of a graph \mathcal{G} is the assignment of a direction to each edge.

in a totally distributed way. In the case of linear coupling $f_{max} = \infty$, the critical value $K_U = 0$ and the network always converges (provided that it is connected).

3. ENERGY CONSUMPTION

All the consensus techniques presented in the previous section have a convergence rate proportional to the algebraic connectivity $\lambda_2(\mathbf{L}_a)$. Hence, from the point of view of convergence time, they can all be treated using the same formulation. In the case of a weighted Laplacian, $\mathbf{L}_a := \mathbf{B}\mathbf{D}_a\mathbf{B}^T$, the weights a_i , could be chosen to maximize the algebraic connectivity $\lambda_2(\mathbf{L}_a)$ of the weighted graph [5]. However, sometimes, as in the case studied in this paper, the weights are simply the channel coefficients and so they are given.

Let us evaluate now the total energy spent to reach the final estimate, within a specified accuracy. For mathematical tractability reasons, we consider the simple case where all the nodes transmit with the same power p_T . In a wireless network, the connectivity (and thus the Laplacian of the graph) depends on the transmit power of each node (as well as the properties of the propagation medium, such as the power-law attenuation factor, fading and shadowing), so that the algebraic connectivity $\lambda_2(\mathbf{L}_a)$ depends on p_T . The convergence time is proportional to $1/\lambda_2(\mathbf{L}_a)$. The total energy consumption is then

$$\mathcal{E} = \frac{N p_T}{\lambda(\mathbf{L}_a(p_T))}, \quad (8)$$

where we have made explicit the dependence of $\lambda(\mathbf{L}_a(p_T))$ on the transmit power p_T .

Let us now see how the network topology changes with p_T . The topological model depends on two sources of randomness, in general: the spatial distribution of the nodes and the channel fading. Furthermore, a link between two nodes is established if the received power is greater than a threshold p_{Rmin} (the SINR model).

Let us consider the case with no fading. The received power p_R depends on the transmit power and on the transmit/receive distance r as $p_R = \frac{p_T}{r^\eta}$, where η is the path loss exponent (typically between 2 and 6). This means that, in the absence of fading, there is a link between any two nodes i and j if their distance r_{ij} is smaller than a critical range $r_0 := \left(\frac{p_T}{p_{Rmin}}\right)^{1/\eta}$. If the nodes are randomly distributed on a plane and the links are established according to the previous rule, the graph describing the network is a *random geometric graph*. We consider a ring topology first, where it is possible to obtain closed form expressions. Then we will consider the more practical planar case.

Ring topology

A ring topology is composed of nodes uniformly spaced over a ring of radius R . This is a regular topology where all the nodes have the same degree, say d . Furthermore, the

Laplacian has a circulant Toeplitz structure that facilitates closed-form computation of the algebraic connectivity:

$$\lambda_2 = 4 \sum_{i=1}^{d/2} \sin^2(\pi i / N) = d + 2 \frac{\sin(\frac{\pi d}{2N})}{\sin(\frac{\pi}{N})} \cos(\frac{\pi}{N}(1 + \frac{d}{2}))$$

For fixed d , λ_2 decreases as N increases; for fixed d/N , it increases with N . In our set-up the degree d is a function of the transmit power, i.e. $d = d(p_T)$. The total energy consumption is then

$$\mathcal{E} = \frac{N p_T}{4 \sum_{i=1}^{d(p_T)/2} \sin^2(\pi i / N)} \quad (9)$$

As an example, in Fig. 1, we report the energy consumption, as a function of the transmit power p_T , for different values of the path loss exponent. Interestingly, we observe that when

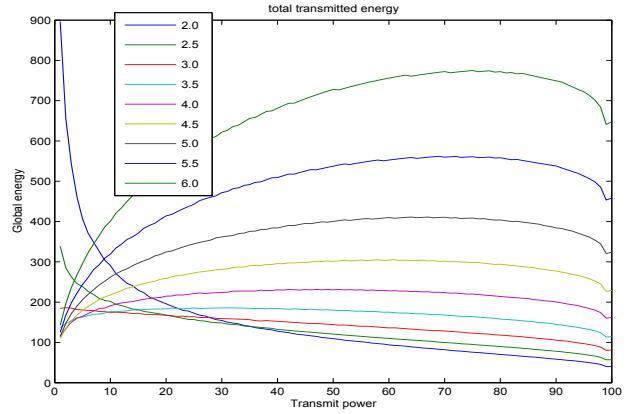


Fig. 1. Total energy consumption as a function of the local transmit power for a ring topology network.

the cost of communication is low (the path exponent is small), it is better to have full connectivity to reach consensus more rapidly. Conversely, when the cost of communication is high (the path loss exponent is high), it is better to have minimum connectivity to save energy.

Planar topology

An interesting question arising in the design of sensor networks is: Given a certain coverage area and number of nodes, is it better to distribute the nodes randomly or over a regular grid? To provide an answer to this question, we considered two planar networks, composed of N nodes distributed over a square of unit side: one with nodes randomly distributed over the square and the other with nodes placed over a rectangular grid. To get rid of any undesired border effect, we consider a toroidal surface. In both cases, we consider a power decaying law $p_R = p_T/(1 + r^2)$, to avoid the unrealistic situation in which the receive power is greater than the transmit power. In Fig. 2, we report the total energy consumption \mathcal{E} , as computed in (8), vs. the transmit power p_T . The number of nodes

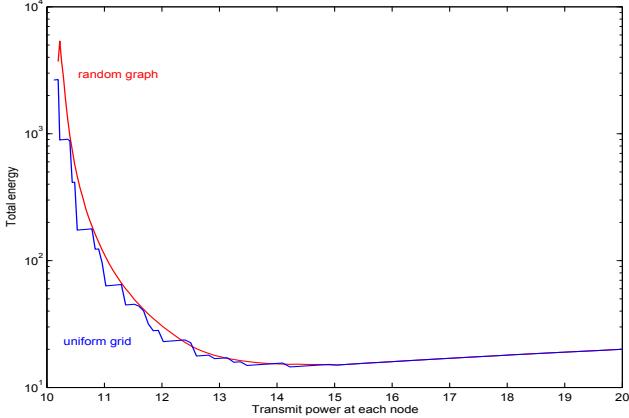


Fig. 2. Total energy consumption as a function of the local transmit power for regular and random networks.

is $N = 100$. In the random graph case, we report a few simulations run over 100 independent generations of the spatial points. From Fig. 2, we can draw two important conclusions: 1) there always exists an optimal p_T that minimizes the overall energy consumption; 2) the random graph requires, practically, the same global energy than the regular graph. As

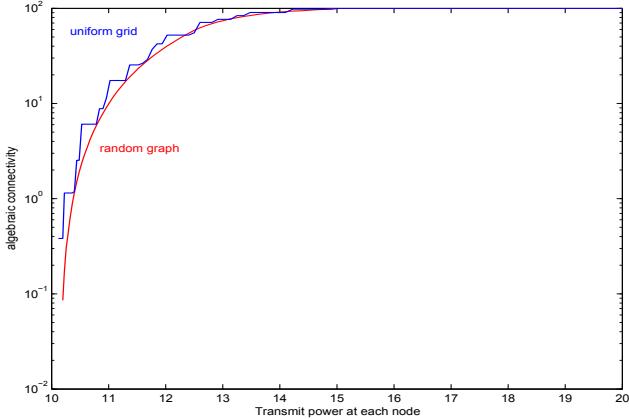


Fig. 3. Algebraic connectivity of the networks of Fig. 1.

expected, the algebraic connectivity for the two topologies is very similar, and this explains the result of Fig. 2.

The behavior shown in Fig. 2 can be explained by recalling the results of [9], where it was shown that all nodes of a random geometric graph, with nodes located over a unit area toroidal surface, tend to have, for large N , the same degree $d = \pi N r_0^2$, with high probability (i.e., with probability greater than $1 - 1/N^2$). But a uniform grid, located again on a unit area toroidal surface, is also a regular graph with degree approximately equal to $\pi N r_0^2$, for large N . Hence, for large N , uniform and random geometric graphs tend to behave similarly. This statement is also confirmed by Fig. 3, where we report the algebraic connectivity for the same

case analyzed in Fig. 2. The results shown in this paper represent only a first step in the minimization of the energy necessary to reach a consensus through a decentralized, iterative, mechanism. Several extensions are worth of being analyzed. We have considered, for simplicity, only the case in which the transmit power is the same for all the nodes. This has been a simplifying assumption that has reduced the number of unknowns to one. In general, it will be interesting to formulate the optimization problem where we minimize the total energy consumption (8) with the respect to the vector of powers transmitted by each node $\mathbf{p} = (p_1, p_2, \dots, p_N)$. Another extension concerns the inclusion of proper channel fading models. Moreover, in sensor networks, energy spent for processing by the receiver can be a significant factor that should properly be taken into account.

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